Patent Claims

1. Compounds of general formula (I),

wherein

or

 R^1 , R^2 , R^{10} , R^{11} independently of one another denote a group selected from among hydrogen, halogen, CN, NO_2 , and -NHCXNH₂ or a group selected from among optionally substituted -COR⁷, -COOR⁷, -CONR⁷R¹³, -OR¹⁴, $NR^{13}R^{15}$, C_1 - C_{10} -alkyl, C_3 - C_8 -cycloalkyl, - $NR^{16}CX$ - R^{17} , - $NR^{18}CX$ - OR^{19} , - $NR^{20}SO_mR^{21}$, - $SO_pNR^{22}R^{23}$ and - SO_qR^{24} .

m, p, q denotes 0, 1 or 2 n denotes 0, 1, 2 or 3

 R^3 denotes hydrogen or a group selected from among optionally substituted C_1 - C_{10} -alkyl, C_6 - C_{10} -aryl, heterocyclyl and C_3 - C_8 -cycloalkyl, -CX- C_1 - C_{10} -alkyl, -CX- C_6 - C_{14} -aryl,

 R^4 , R^5 independently of one another denote hydrogen, halogen or optionally substituted C_1 - C_{10} -alkyl,

R⁴ and R⁵ together denote a C₃-C₈-alkyl bridge,

R⁶ denotes a group selected from among the general formulae

l,k independently of one another denote 1,2 or 3,

R²⁵, R²⁶, R²⁷, R²⁸ independently of one another denote a group selected from among hydrogen, OH, halogen, CN and NO₂, or

a group selected from among optionally substituted C_1 - C_{10} -alkyl, C_6 - C_{18} -aryl, heteroaryl, heterocyclyl, -CX- R^{17} , -OR¹⁴, NR¹³R¹⁵, C_2 - C_8 -cycloalkyl, -NR²⁰SO_mR²¹, -SO_pNR²²R²³, -SO_qR²⁴, -NR¹⁸CX- R^{19} , -NR¹⁸CXOR¹⁷, while R^{25} and R^{26} cannot simultaneously denote hydrogen,

 R^8 denotes hydrogen or a group selected from among optionally substituted C₁-C₁₀-alkyl, C₆-C₁₈-aryl, -SO_q- C₁-C₁₀-alkyl, -SO_q-C₆-C₁₄-aryl, -CX- C₁-C₁₀-alkyl, -CX-C₆-C₁₄-aryl, C₆-C₁₀-aryl, heterocyclyl and C₃-C₈-cycloalkyl

 R^9 denotes hydrogen or a group selected from among optionally substituted C_1 - C_{10} -alkyl, C_6 - C_{14} -aryl, heteroaryl, C_3 - C_8 -cycloalkyl and heterocycloalkyl,

 R^{12} denotes hydrogen or a group selected from among optionally substituted benzyl, C_1 - C_{12} -alkyl and C_6 - C_{14} -aryl,

 R^7 , R^{13} , R^{15} , R^{16} , R^{18} , R^{20} . R^{22} , R^{23} independently of one another denote hydrogen, or a group selected from among optionally substituted C_1 - C_{10} -alkyl, C_6 - C_{14} -aryl, heterocyclyl and C_3 - C_8 -cycloalkyl

 R^{14} , R^{19} , R^{29} independently of one another denote hydrogen or a group selected from among optionally substituted C_1 - C_{10} -alkyl, C_6 - C_{14} -aryl, C_3 - C_8 -cycloalkyl, heteroaryl, heterocyclyl, -CXNR₁₃R₁₅ and -CXR₇

 R^{17} denotes a group selected from among C_1 - C_{10} -alkyl, C_6 - C_{14} -aryl, heterocyclyl, heteroaryl and C_3 - C_8 -cycloalkyl

 R^{21} , R^{24} independently denote hydrogen or OH, or a group selected from among optionally substituted $N(C_1-C_{10}-alkyl)_2$, $N(C_3-C_8-cycloalkyl)$, $C_1-C_{10}-alkyl$, $C_6-C_{14}-aryl$, heterocyclyl, heteroaryl and $C_3-C_8-cycloalkyl$

and

X denotes O, S or NR²⁹,

optionally in the form of the tautomers, the racemates, the enantiomers, the diastereomers and the mixtures thereof, as well as optionally the pharmacologically acceptable acid addition salts thereof.

2. Compounds according to claim 1, wherein

R^{10,} R¹¹ independently of one another denote hydrogen or halogen,

m, p, q independently of one another denote 0, 1 or 2

n denotes 0, 1, 2 or 3

R³ denotes hydrogen or C₁-C₅-alkyl

R⁴, R⁵ independently of one another denote hydrogen or C₁-C₅-alkyl,

 R^8 denotes a group selected from among hydrogen, C_1 - C_5 -alkyl, - SO_q - C_1 - C_5 -alkyl, - SO_q - C_6 - C_{14} -aryl, phenyl and C_3 - C_6 -cycloalkyl

R⁹ denotes hydrogen or C₁-C₁₀-alkyl

R¹² denotes hydrogen or benzyl

R¹³, R¹⁵, R¹⁶, R¹⁸ independently of one another denote a group selected from among hydrogen, C₁-C₅-alkyl, C₃-C₆-cycloalkyl and phenyl

 R^{14} , R^{19} independently of one another denote hydrogen or C_1 - C_5 -alkyl,

and

 R^{17} denotes optionally substituted C_1 - C_5 -alkyl or C_6 - C_{10} -aryl.

3. Compounds according to claim 1 or 2, wherein

R¹⁰, R¹¹ denote hydrogen

m, p, q denote 0, 1 or 2

n denotes 0, 1, 2 or 3

R³ denotes hydrogen

R⁴, R⁵ independently of one another denote hydrogen or methyl,

R⁸ denotes hydrogen, -SO_q-C₆-C₁₄-aryl or -SO₂-C₁-C₅-alkyl

R⁹ denotes hydrogen

R¹² denotes hydrogen or benzyl,

R¹³, R¹⁵, R¹⁶, R¹⁸ independently of one another denote a group selected from among hydrogen, C₁-C₁₅-alkyl and phenyl,

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R¹⁴, R¹⁹ independently of one another denote hydrogen or C₁-C₅-alkyl,

and

 R^{17} denotes C_1 - C_5 -alkyl or C_6 - C_{14} -aryl.

4. Compounds according to one of claims 1 to 3, wherein

R¹ denotes a group selected from among hydrogen, NO₂, NH₂, -NHCX-R¹⁷ and -NHSO₂R²¹

R² denotes hydrogen or halogen

n denotes 2,

R³ denotes hydrogen

R⁴, R⁵ denote hydrogen or methyl

R⁶ denotes a group selected from among the general formulae

I,k denote 1

R²⁶, R²⁷ denote hydrogen,

R⁸ denotes hydrogen or -SO₂CH₃,

R⁹ denotes hydrogen,

R^{10,} R¹¹ denote hydrogen, and

R¹² denotes hydrogen or benzyl.

- 5. Compounds according to one of claims 1 to 4, wherein
- R⁶ denotes a group selected from among the general formulae

- 6. Compounds according to one of claims 1 to 5, wherein
- R⁶ denotes an optionally substituted group of formula (j)

- 7. Compounds of formula (I) according to one of claims 1 to 6 for use as pharmaceutical compositions.
- 8. Compounds of formula (I) according to one of claims 1 to 6 for use as pharmaceutical compositions with a selective beta-3-agonistic activity.
- 9. Use of a compound of formula (I) according to one of claims 1 to 6 for preparing a pharmaceutical composition for the treatment and/or prevention of diseases connected with the stimulation of beta-3-receptors.
- 10. Method for the treatment and/or prevention of diseases connected with the stimulation of beta-3-receptors, characterised in that an effective amount

of a compound of formula I according to claim 1 to 6 is administered to a patient.

- 11. Pharmaceutical composition, containing as active substance one or more compounds of general formula (I) according to one of claims 1 to 6 or the physiologically acceptable salts thereof optionally combined with conventional excipients and/or carriers.
- 12. Pharmaceutical composition containing as active substance one or more compounds of general formula (I) according to one of claims 1 to 6 or the physiologically acceptable salts thereof and one or more active substances selected from among antidiabetics, inhibitors of protein tyrosinephosphatase 1, substances which influence deregulated glucose production in the liver, lipid lowering agents, cholesterol absorption inhibitors, HDL-raising compounds, active substances for the treatment of obesity and modulators or stimulators of the adrenergic via alpha 1 and alpha 2 as well as beta 1, beta 2 and beta 3 receptors.
- 13. Process for preparing a compound of general formula (I),

$$R^{12}$$
 R^{10}
 R

wherein

R¹-R²⁸ and X may have the meanings given in claims 1 to 6, characterised in that a compound of general formula (II)

wherein

R⁴ and R⁵ have the meanings given in claims 1 to 6, is converted by means of a chlorinating agent into a compound of formula (III)

$$R^8$$
 R^4
 R^5
(III)

the compound of formula (III), optionally provided with an amino protective group, is reacted with an optionally substituted compound selected from among the general formulae (IVa) to (IVi)

$$(IV) \\ H \sim (R^{28})_{k} \qquad G$$

$$(R^{25})_{i} \qquad H \sim (R^{28})_{k} \qquad G$$

$$(R^{27})_{i} \qquad H \sim (R^{28})_{k} \qquad H \sim (R^{28})_{k} \qquad H \sim (R^{28})_{k} \qquad G$$

$$(R^{27})_{i} \qquad H \sim (R^{28})_{k} \qquad H \sim (R^{28})_{k} \qquad G$$

$$(R^{27})_{i} \qquad G$$

$$(R^{27})_{i} \qquad G$$

wherein

k, I, R^{27} and R^{28} have the meanings given in claims 1 to 6, and the product of formula (V)

$$\begin{array}{c} \text{(V)} \\ \text{H-N} \\ \text{R}^4 \\ \text{R}^5 \end{array}$$

wherein n, R^4 , R^5 , R^6 and R^8 have the meanings given in claims 1 to 6, is reacted with a compound of formula (VI)

$$R^{12} \longrightarrow R^{10}$$

$$R^{12} \longrightarrow R^{10}$$

$$R^{12} \longrightarrow R^{2}$$

wherein R¹, R², R⁹ and R¹⁰ to R¹² have the meanings given in claims 1 to 6.